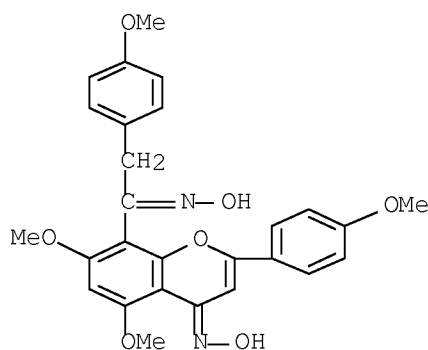


ACCESSION NUMBER: 1956:89181 CAPLUS Full-text  
 DOCUMENT NUMBER: 50:89181  
 ORIGINAL REFERENCE NO.: 50:16759f-g  
 TITLE: Structure of sciadopitysin, a flavonoid from the leaves of *Sciadopitys verticillata*. III. The structure of oxoflavone and carboxyflavone  
 AUTHOR(S): Kariyone, Tatsuo; Kawano, Nobusuke  
 SOURCE: Yakugaku Zasshi (1956), 76, 453-6  
 CODEN: YKKZAJ; ISSN: 0031-6903  
 DOCUMENT TYPE: Journal  
 LANGUAGE: Unavailable  
 AB IV and V were each shown to have 2 MeO groups and 2 phenolic HO groups and were assumed to have the skeleton of acacetin 7-Me ether (VI) with an acyl group as the side-chain. IV is VI where the acyl is COCH<sub>2</sub>C<sub>6</sub>H<sub>4</sub>OH and V is IV bearing a group on the VI moiety.  
 IT 874530-27-5P, Flavone, 4',5,7-trimethoxy-8-[(p-methoxyphenyl)acetyl]-, dioxime  
 RL: PREP (Preparation)  
 (preparation of)  
 RN 874530-27-5 CAPLUS  
 CN Flavone, 4',5,7-trimethoxy-8-[(p-methoxyphenyl)acetyl]-, dioxime (5CI)  
 (CA INDEX NAME)

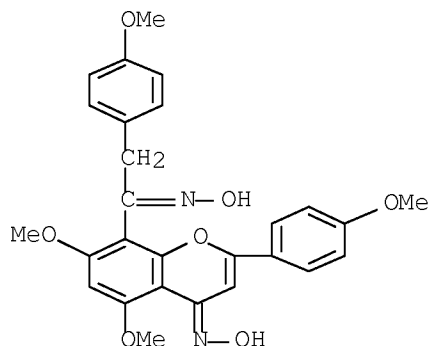


L9 ANSWER 55 OF 66 CAPLUS COPYRIGHT 2007 ACS on STN  
 ACCESSION NUMBER: 1956:89180 CAPLUS Full-text  
 DOCUMENT NUMBER: 50:89180  
 ORIGINAL REFERENCE NO.: 50:16759e-f  
 TITLE: Structure of sciadopitysin, a flavonoid from the leaves of *Sciadopitys verticillata*. II. Degradation of sciadopitysin  
 AUTHOR(S): Kariyone, Tatsuo; Kawano, Nobusuke  
 SOURCE: Yakugaku Zasshi (1956), 76, 451-2  
 CODEN: YKKZAJ; ISSN: 0031-6903  
 DOCUMENT TYPE: Journal  
 LANGUAGE: Unavailable  
 AB Decomposition of I by boiling 1.5 hrs. with 20% aqueous KOH afforded an oxoflavone (IV), C<sub>25</sub>H<sub>20</sub>O<sub>7</sub>, yellow columns, m. 241-2°, a carboxyflavone (V), C<sub>26</sub>H<sub>20</sub>O<sub>9</sub>, yellow, m. 311° (decomposition), anisic acid, p-MeOC<sub>6</sub>H<sub>4</sub>Ac, and 4,2,6-MeO-(HO)<sub>2</sub>C<sub>6</sub>H<sub>2</sub>Ac; the yield of these products varied with the concentration of KOH and duration of boiling.  
 IT 874530-27-5P, Flavone, 4',5,7-trimethoxy-8-[(p-methoxyphenyl)acetyl]-, dioxime

RL: PREP (Preparation)  
(preparation of)

RN 874530-27-5 CAPLUS

CN Flavone, 4',5,7-trimethoxy-8-[(p-methoxyphenyl)acetyl]-, dioxime (5CI)  
(CA INDEX NAME)



L9 ANSWER 56 OF 66 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1956:89179 CAPLUS Full-text

DOCUMENT NUMBER: 50:89179

ORIGINAL REFERENCE NO.: 50:16759d-f

TITLE: Structure of sciadopitysin, a flavonoid from the  
leaves of *Sciadopitys verticillata*. I

AUTHOR(S): Kariyone, Tatsuo; Kawano, Nobusuke

SOURCE: Yakugaku Zasshi (1956), 76, 448-50

CODEN: YKKZAJ; ISSN: 0031-6903

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

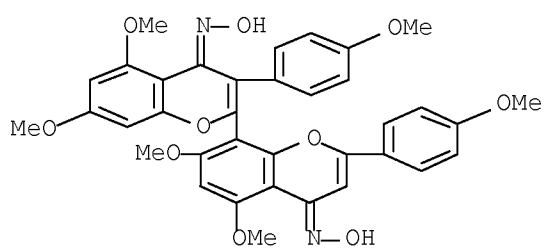
AB Air-dried leaves (3 kg.) in 20 l. CHCl<sub>3</sub>:CCl<sub>2</sub> refluxed 3 hrs., the extract concentrated to 3 l., allowed to stand overnight, the waxy precipitate warmed up, and the insol. residue filtered off gave 9 g. sciadopitysin (I), yellow columns, m. 285-6° (from C<sub>5</sub>H<sub>5</sub>NEtOH); triacetate, prisms, m. 264°; tri-Me ether (IIIa), m. 214-15° (oxime, C<sub>36</sub>H<sub>32</sub>O<sub>10</sub>N<sub>2</sub>, columns, m. 248-9°). I (0.1 g.) in 40 ml. Me<sub>2</sub>CO treated with 0.5 g. each of K<sub>2</sub>CO<sub>3</sub> and MeI and the mixture refluxed 40 min. and concentrated gave a I mono-Me ether (II), m. 282°, identical with ginkgetin di-Me ether (III) by mixed m.p. II gives a diacetate, m. 228°, identical with the diacetate of III by mixed m.p. Demethylation of I with HI gave C<sub>30</sub>H<sub>18</sub>O<sub>10</sub>.5H<sub>2</sub>O, m. above 360°, and acetylation of this substance gave a product, m. 240°, identical with that prepared by treating ginkgetin in a similar way.

IT 854209-81-7P, [2,8'-Bi-4H-1-benzopyran]-4,4'-dione, 5,5',7,7'-tetramethoxy-2',3-bis(p-methoxyphenyl)-, dioxime  
874530-27-5P, Flavone, 4',5,7-trimethoxy-8-[(p-methoxyphenyl)acetyl]-, dioxime

RL: PREP (Preparation)  
(preparation of)

RN 854209-81-7 CAPLUS

CN [2,8'-Bi-4H-1-benzopyran]-4,4'-dione, 5,5',7,7'-tetramethoxy-2',3-bis(p-methoxyphenyl)-, dioxime (5CI) (CA INDEX NAME)



RN 874530-27-5 CAPLUS

CN Flavone, 4',5,7-trimethoxy-8-[(p-methoxyphenyl)acetyl]-, dioxime (5CI)  
(CA INDEX NAME)

